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                status data
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        MAY 28 CAS databases on STN enhanced with NANO super role in
                records back to 1992
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NEWS EXPRESS MAY 26 09 CURRENT WINDOWS VERSION IS V8.4, AND CURRENT DISCOVER FILE IS DATED 06 APRIL 2009.

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=> file reg

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0.22

0.22

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STRUCTURE FILE UPDATES: 15 JUN 2009 HIGHEST RN 1158168-92-3 DICTIONARY FILE UPDATES: 15 JUN 2009 HIGHEST RN 1158168-92-3

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=>

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chain nodes :

10 11 12 14 22 26 27 28 31 33 35

ring nodes :

1 2 3 4 5 6 7 8 9 13 16 17 18 19 20

chain bonds :

ring bonds :

exact/norm bonds :

 $4-26 \quad 5-7 \quad 6-9 \quad 7-8 \quad 7-10 \quad 8-9 \quad 8-11 \quad 9-33 \quad 9-35 \quad 10-22 \quad 12-14 \quad 27-28$

exact bonds : 11-12 12-13

normalized bonds :

 $1-2 \quad 1-6 \quad 2-3 \quad 3-4 \quad 4-5 \quad 5-6 \quad 13-16 \quad 13-20 \quad 16-17 \quad 17-18 \quad 18-19 \quad 19-20$

G1:H,CH3,Et

G2:H,CH3,Et,X

G3:H,[*1]

Connectivity:

14:1 E exact RC ring/chain

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS 11:CLASS 12:CLASS 13:Atom 14:CLASS 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom

22:CLASS 26:CLASS

Generic attributes :

28:

Saturation : Saturated

L1 STRUCTURE UPLOADED

=> d

L1 HAS NO ANSWERS

L1 STR

G1 H, Me, Et

G2 H, Me, Et, X

G3 H, [@1]

Structure attributes must be viewed using STN Express query preparation.

=> s 11

SAMPLE SEARCH INITIATED 10:41:51 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 136 TO ITERATE

100.0% PROCESSED 136 ITERATIONS

INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

2021 TO 3419 PROJECTED ITERATIONS: PROJECTED ANSWERS: 1882 TO 3238

50 SEA SSS SAM L1 L2

=> s 11 full

FULL SEARCH INITIATED 10:42:17 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 3010 TO ITERATE

100.0% PROCESSED 3010 ITERATIONS 2863 ANSWERS

SEARCH TIME: 00.00.01

L3 2863 SEA SSS FUL L1

=> save temp 13

ENTER NAME OR (END):yc10590976/a

ANSWER SET L3 HAS BEEN SAVED AS 'YC10590976/A'

=> file caplus

COST IN U.S. DOLLARS

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187.06

186.84

50 ANSWERS

FULL ESTIMATED COST

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FILE COVERS 1907 - 16 Jun 2009 VOL 150 ISS 25 FILE LAST UPDATED: 15 Jun 2009 (20090615/ED) REVISED CLASS FIELDS (/NCL) LAST RELOADED: Feb 2009 USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Feb 2009

CAplus now includes complete International Patent Classification (IPC) reclassification data for the third quarter of 2008.

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s 13 26 L3 L4=> s halide salt 163353 HALIDE 135680 HALIDES 235475 HALIDE (HALIDE OR HALIDES) 890769 SALT 676873 SALTS 1314616 SALT (SALT OR SALTS) L52345 HALIDE SALT (HALIDE(W)SALT) \Rightarrow s 14 and 15 1 L4 AND L5 => s 14 and electrolyte 288913 ELECTROLYTE 148781 ELECTROLYTES 345301 ELECTROLYTE (ELECTROLYTE OR ELECTROLYTES) L7 1 L4 AND ELECTROLYTE => s 16 not 17 0 L6 NOT L7 => s 14 and decomposition 209589 DECOMPOSITION 1263 DECOMPOSITIONS 210458 DECOMPOSITION (DECOMPOSITION OR DECOMPOSITIONS) 466936 DECOMPN 5163 DECOMPNS 468728 DECOMPN (DECOMPN OR DECOMPNS) 559969 DECOMPOSITION (DECOMPOSITION OR DECOMPN) L91 L4 AND DECOMPOSITION => s 19 not 17 L10 0 L9 NOT L7 => d 16 isib abs hitstr tot 'ISIB' IS NOT A VALID FORMAT FOR FILE 'CAPLUS' The following are valid formats: ABS ----- GI and AB ALL ----- BIB, AB, IND, RE

APPS ----- AI, PRAI

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IND ----- Indexing data
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=> d 16 ibib abs hitstr tot

ACCESSION NUMBER: 2005:1004569 CAPLUS <u>Full-text</u>

DOCUMENT NUMBER: 143:292577

TITLE: Composition containing benzamidine derivative and

method for stabilizing benzamidine derivative

į

INVENTOR(S): Suzuki, Yasuyuki; Fujioka, Satoshi

PATENT ASSIGNEE(S): Eisai Co., Ltd., Japan SOURCE: PCT Int. Appl., 29 pp.

CODEN: PIXXD2

DOCUMENT TYPE: Patent LANGUAGE: Japanese

FAMILY ACC. NUM. COUNT: 1 **Current application**

PATENT INFORMATION:

PA:	PATENT NO.					KIND DATE		APPLICATION NO.						DATE				
						WO 2005-JP3742						20050304						
	W:	ΑE,	AG,	AL,	AM,	AT,	ΑU,	ΑZ,	BA,	BB,	BG,	BR,	BW,	ΒY,	BZ,	CA,	CH,	
		CN,	CO,	CR,	CU,	CZ,	DE,	DK,	DM,	DZ,	EC,	EE,	EG,	έs,	FI,	GB,	GD,	
		GE,	GH,	GM,	HR,	HU,	ID,	IL,	IN,	IS,	JP,	ΚE,	KG,	KΡ,	KR,	KΖ,	LC,	
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		NO,	NZ,	OM,	PG,	PH,	PL,	PT,	RO,	RU,	SC,	SD,	SE,	SG,	SK,	SL,	SM,	
		SY,	ТJ,	TM,	TN,	TR,	TT,	TZ,	UA,	UG,	US,	UZ,	VC.	VN,	YU,	ZA,	ZM,	$Z\overline{W}$
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		AZ,	BY,	KG,	KΖ,	MD,	RU,	ТJ,	TM,	ΑT,	BE,	BG,	СĦ,	CY,	CZ,	DE,	DK,	
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AU				A1 20050915				AU 2005-2190∮0					20050304					
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EP	P 1721610 B1 20090513 /																	
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		HR,	LV,	MK,	YU													
CN	CN 1925861			Α				CN 2005-\$0006933					20050304					
	AT 431150			_					AT 2005∜720014					20050304				
US	US 20070208016				A1	A1 20070906			US 2006-590976					20060828				
KR	KR 2006117365			Α		20061116			KR 2006-717795					20060901				
KR	KR 760448			В1		2007	1004											
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										WO 2	005-	JP37	42	,	W 2	0050	304	
THER SO	ER SOURCE(S):				MARPAT 143:292577													

GI

AB Disclosed is a compn. contg. a benzamidine deriv. which is not decompd. even under humidified conditions. Also disclosed is a method for stabilizing a benzamidine deriv. Decompn. reaction of benzamidine derivs. can be suppressed by adding at least one electrolyte selected from the group consisting of

halide salts of alkali metals or alk. earth metals and perchlorates of alkali metals or alk. earth metals to a benzamidine deriv. represented by the general formula I (R1, R2 = H, methoxy, ethoxy; X = H, halogen; Ar = Me, Et, methoxy, ethoxy, tert-Bu, morpholinyl, etc), or a pharmacol. acceptable salt thereof. For example, tablets were prepd. from 1-(3-tert-butyl-4-methoxy-5-morpholino-phenyl)-2- (5,6-diethoxy-7-fluoro-1-imino-1,3-dihydro-isoindol-2-yl)-ethanone 1, lactose 117, hydroxypropyl cellulose 7.5, hydroxypropyl Me cellulose 4.5, NaCl 4.5, cryst. cellulose 15, and magnesium stearate 0.75 g.

IT 474624-49-2 734528-58-6 751475-53-3 752978-65-7 759452-62-5 790653-73-5 792182-11-7

RL: BSU (Biological study, unclassified); THU (Therapeutic use); BIOL (Biological study); USES (Uses)

(compns. contg. benzamidine derivs. and electrolytes, and method for stabilizing benzamidine deriv.)

RN 474624-49-2 CAPLUS

CN Ethanone, 2-(5,6-diethoxy-7-fluoro-1,3-dihydro-1-imino-2H-isoindol-2-yl)-1[3-(1,1-dimethylethyl)-5-[(3S,4S)-3-hydroxy-4-methoxy-1-pyrrolidinyl]-4methoxyphenyl]- (CA INDEX NAME)

Absolute stereochemistry.

RN 734528-58-6 CAPLUS

CN Ethanone, 2-(5,6-diethoxy-7-fluoro-1,3-dihydro-1-imino-2H-isoindol-2-yl)-1[3-(1,1-dimethylethyl)-5-(4-hydroxy-1-piperidinyl)-4-methoxyphenyl]- (CA INDEX NAME)

RN 751475-53-3 CAPLUS

CN Ethanone, 2-(5,6-diethoxy-7-fluoro-1,3-dihydro-1-imino-2H-isoindol-2-yl)-1[3-(1,1-dimethylethyl)-4-methoxy-5-(4-morpholinyl)phenyl]- (CA INDEX NAME)

RN 752978-65-7 CAPLUS

CN Ethanone, 2-(5,6-diethoxy-7-fluoro-1,3-dihydro-1-imino-2H-isoindol-2-yl)-1[3-(1,1-dimethylethyl)-5-[4-(2-hydroxyacetyl)-1-piperazinyl]-4methoxyphenyl]- (CA INDEX NAME)

RN 759452-62-5 CAPLUS

CN Ethanone, 2-(5,6-diethoxy-7-fluoro-1,3-dihydro-1-imino-2H-isoindol-2-yl)-1[3-(1,1-dimethylethyl)-4-methoxy-5-(1-piperazinyl)phenyl]- (CA INDEX NAME)

RN 790653-73-5 CAPLUS

CN 1-Piperazineacetic acid, 4-[5-[2-(5,6-diethoxy-7-fluoro-1,3-dihydro-1-imino-2H-isoindol-2-yl)acetyl]-3-(1,1-dimethylethyl)-2-methoxyphenyl]-, ethyl ester (CA INDEX NAME)

$$\begin{array}{c} \text{EtO} \\ \text{EtO} \\ \end{array} \begin{array}{c} \text{N} \\ \text{CH}_2 \\ \end{array} \begin{array}{c} \text{C} \\ \text{NH} \\ \end{array}$$

RN 792182-11-7 CAPLUS

CN 1-Piperazineacetonitrile, 4-[5-[2-(5,6-diethoxy-7-fluoro-1,3-dihydro-1-imino-2H-isoindol-2-yl)acetyl]-3-(1,1-dimethylethyl)-2-methoxyphenyl]-(CA INDEX NAME)

REFERENCE COUNT: 13 THERE ARE 13 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

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